



# A Strategy to Optimize Local Phase Transformation Strengthening for Next Generation Superalloys

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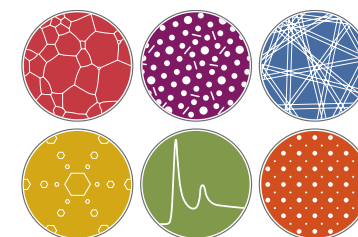
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UNIVERSITY**

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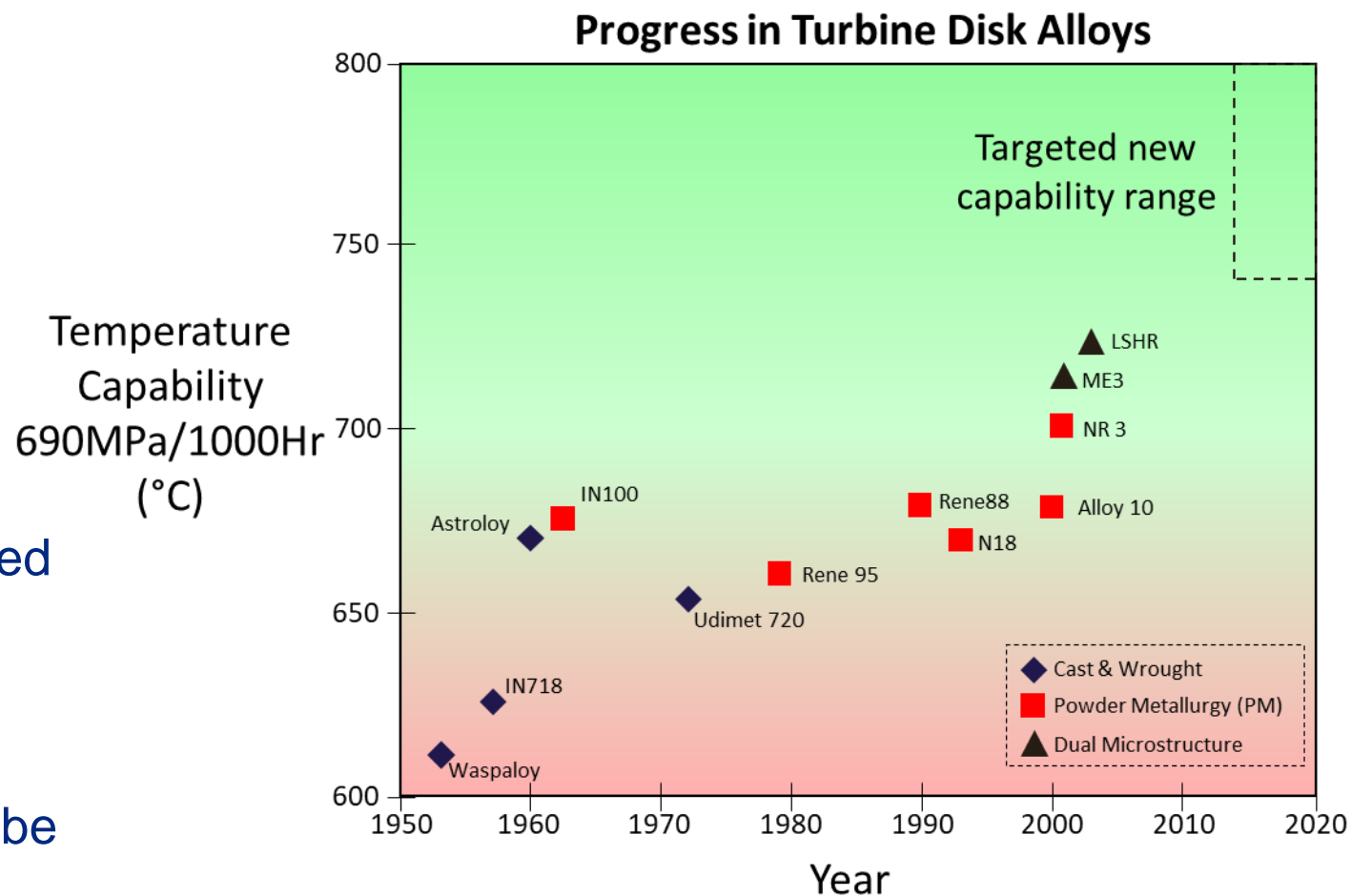
Center for Electron Microscopy and Analysis





# Motivation for Mechanistic Studies

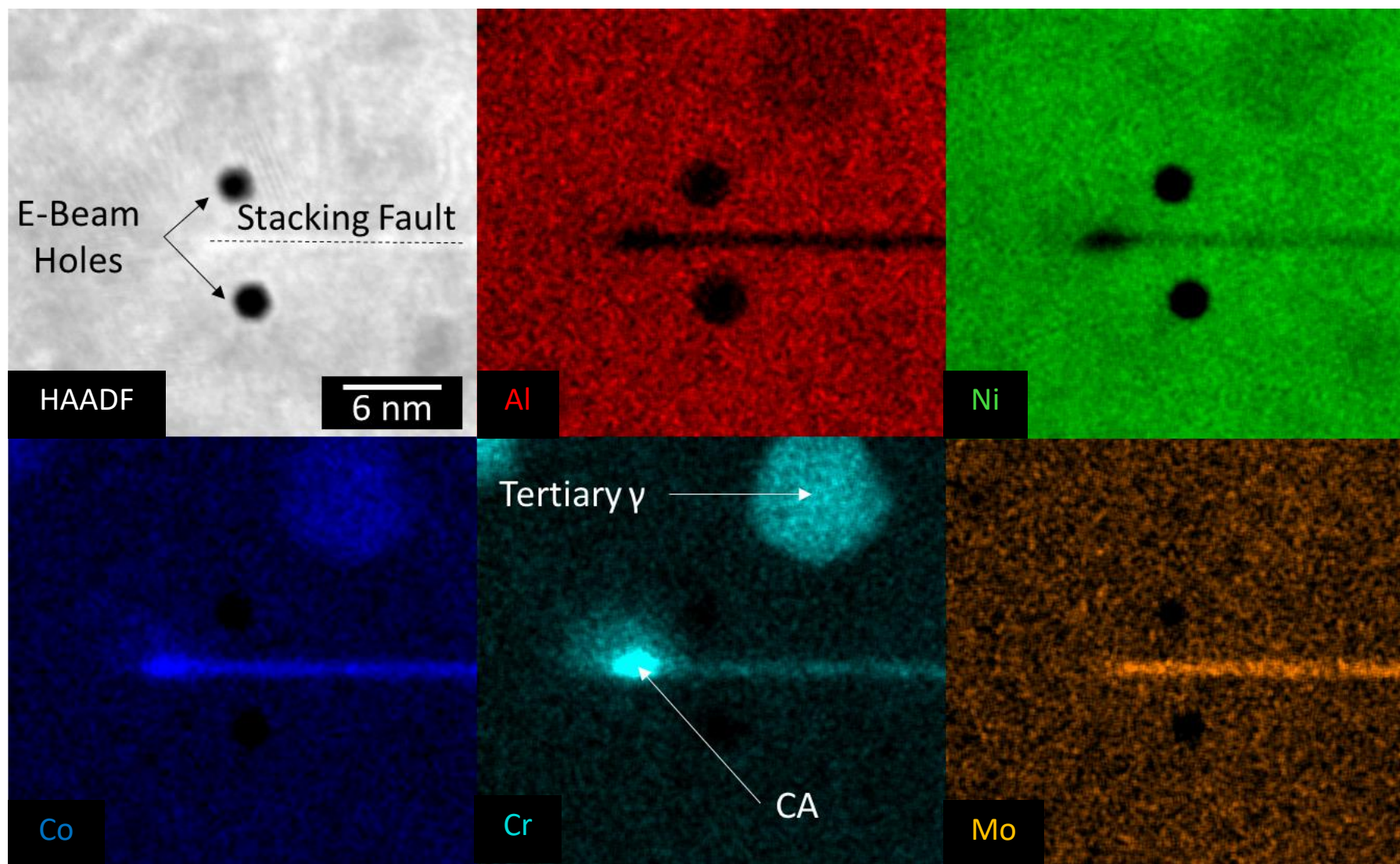
- Material advancements are required to accommodate the higher compressor exit temperatures in jet turbine engines ( $>700^{\circ}\text{C}$  near the rotor rim) for improved efficiency and pollution reduction.
- New deformation mechanisms will become dominant at these higher operating temperatures along with a need for improved creep properties in these disk alloys.
- New understanding and materials will be needed for future advancements





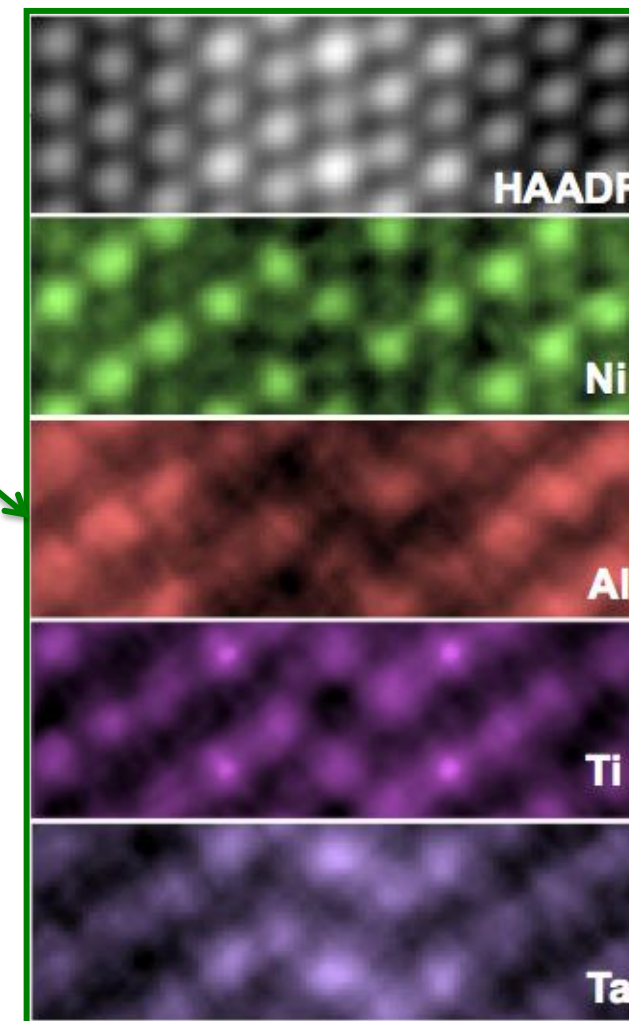
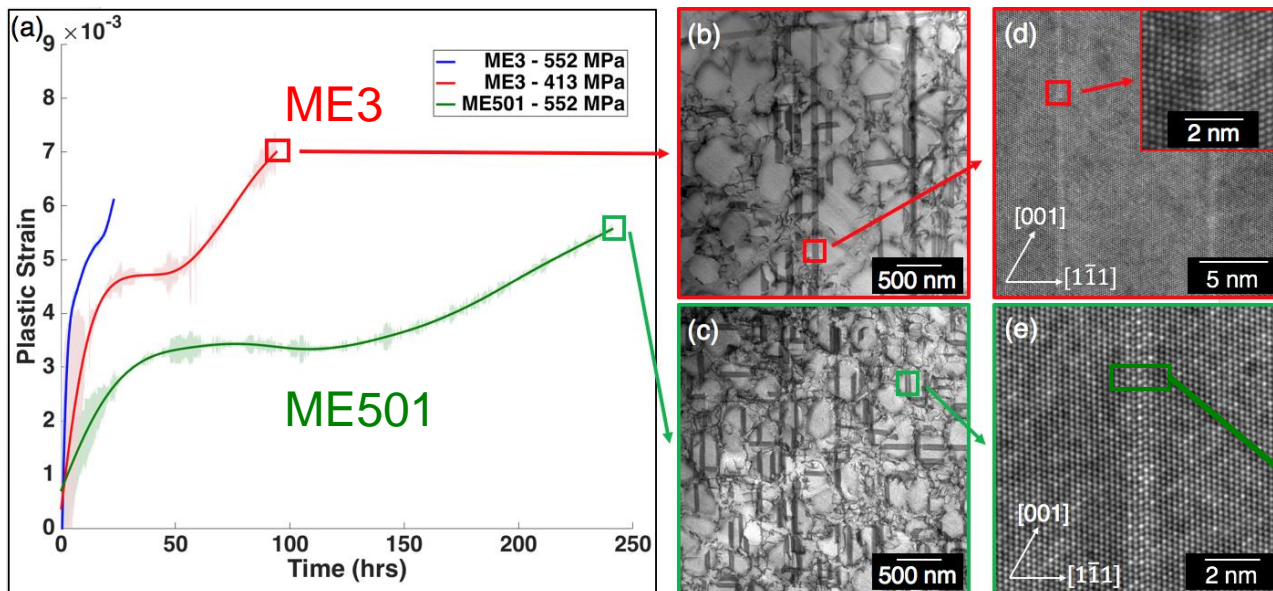
# Segregation along Stacking Faults

Segregation along superlattice stacking faults has been observed in numerous Ni and Co-based superalloys.



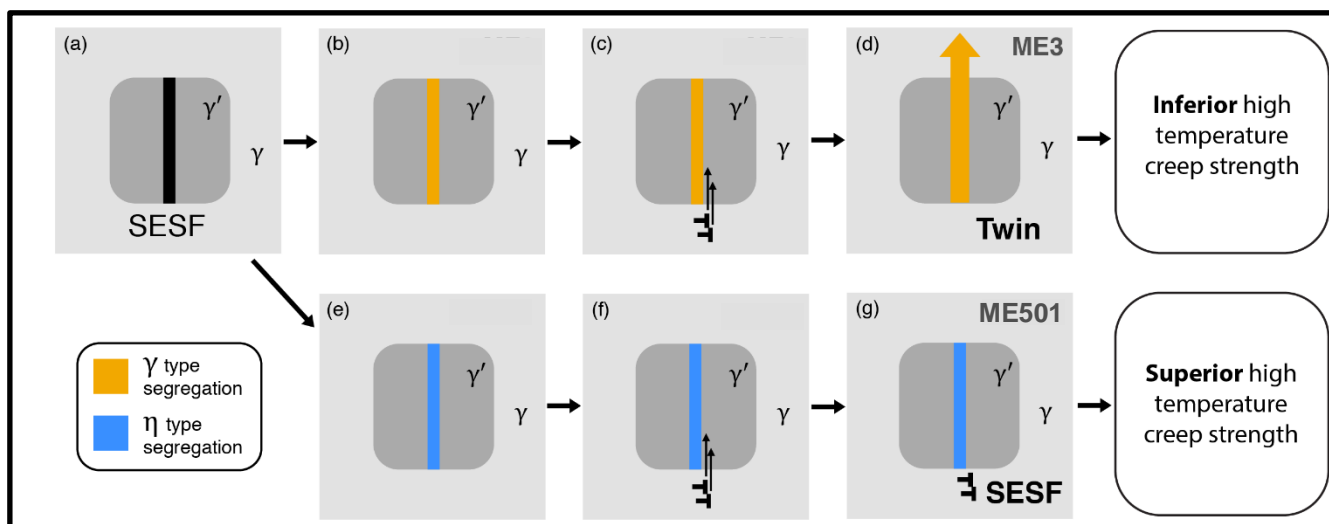


# Phase Transformation Strengthening



**New insight into alloy effects:**

- Segregation of  $\gamma$  formers in ME3 promotes microtwinning
- Formation of  $\eta$  phase at faults in ME501 inhibits microtwinning and improves creep strength

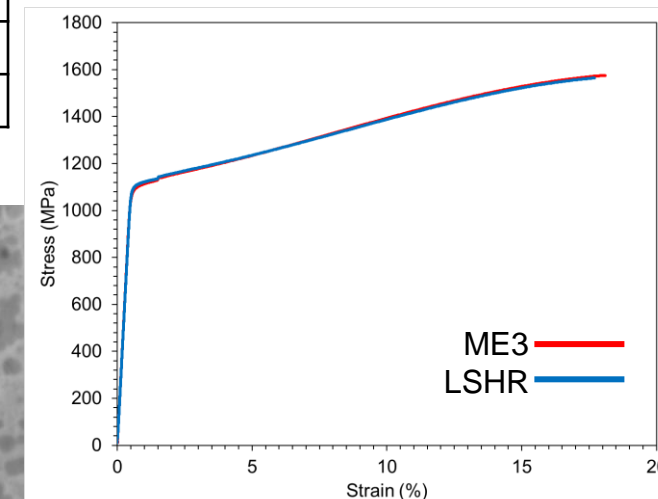




# Material Preparation

Average Alloy Composition in Weight Percent												
Alloy	Cr	Co	Al	Ti	Nb	Mo	Ta	W	Zr	B	C	Ni
LSHR	12.5	20.4	3.5	3.5	1.5	2.7	1.5	4.3	0.05	0.03	0.045	Bal
ME3	13	21	3.4	3.8	0.8	3.7	2.4	2.1	0.05	0.02	0.05	Bal

## Room Temperature Tensile

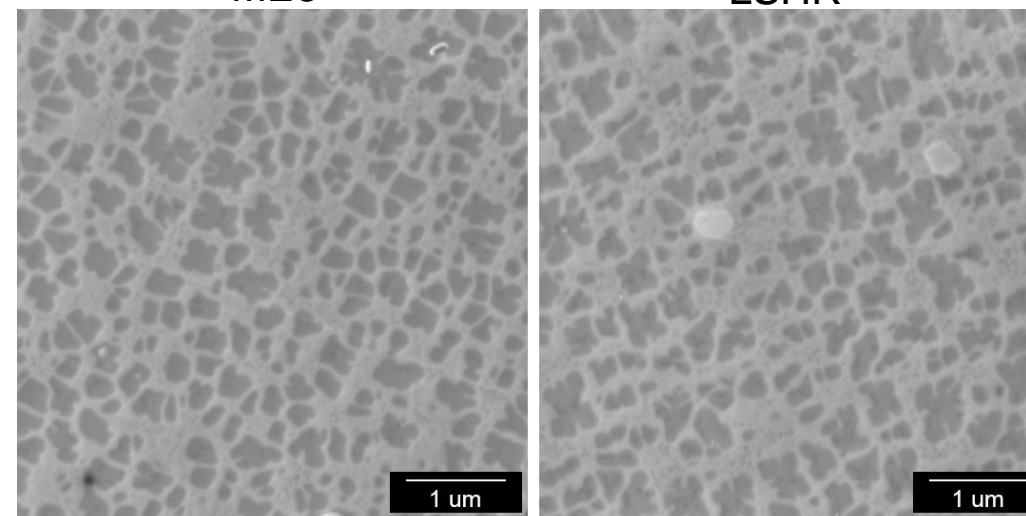
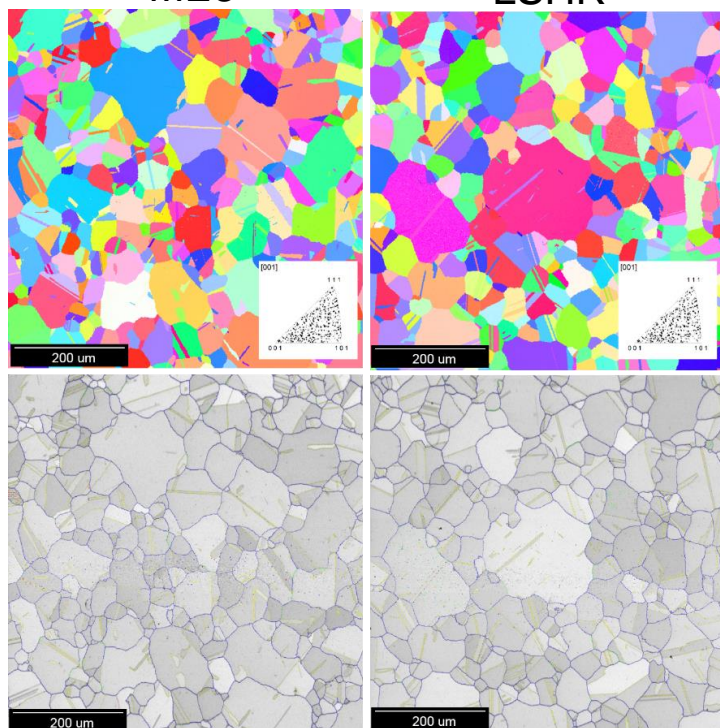


ME3

LSHR

ME3

LSHR



Alloy	Secondary $\gamma'$ VF	Tertiary $\gamma'$ VF	Total $\gamma'$ VF	Average Secondary $\gamma'$ Size	Average Tertiary $\gamma'$ Size
ME3	43.97 ± .6	2.65 ± .4	46.61 ± 1.0	135 nm	15.4 nm
LSHR	43.52 ± 1.7	2.27 ± .1	45.80 ± 1.8	154 nm	15.9 nm

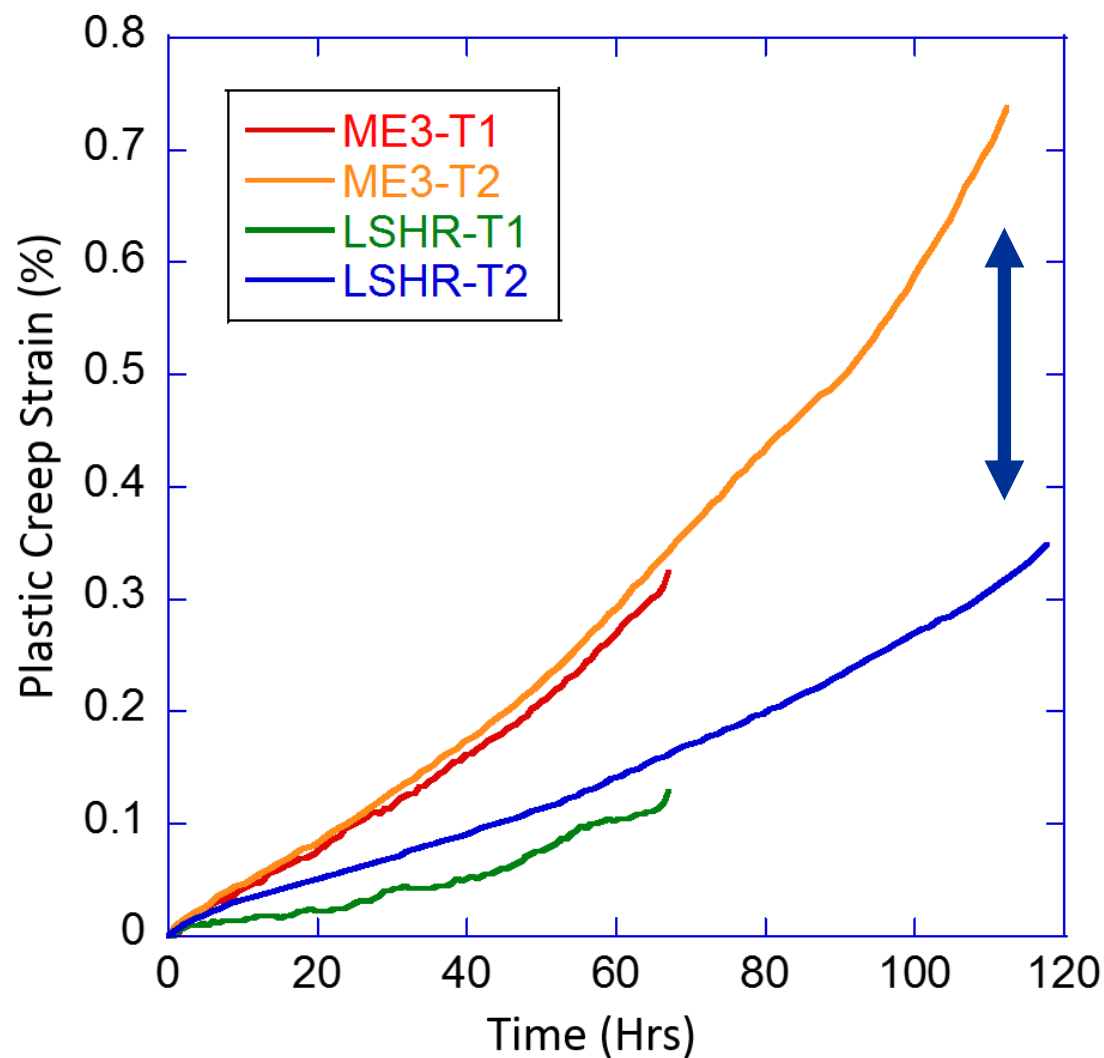
ME3 Average Grain Diameter = 59.2  $\mu\text{m}$   
 LSHR Average Grain Diameter = 59.9  $\mu\text{m}$

The two alloys are microstructurally comparable!

Smith, et al. Acta Materialia, 2019



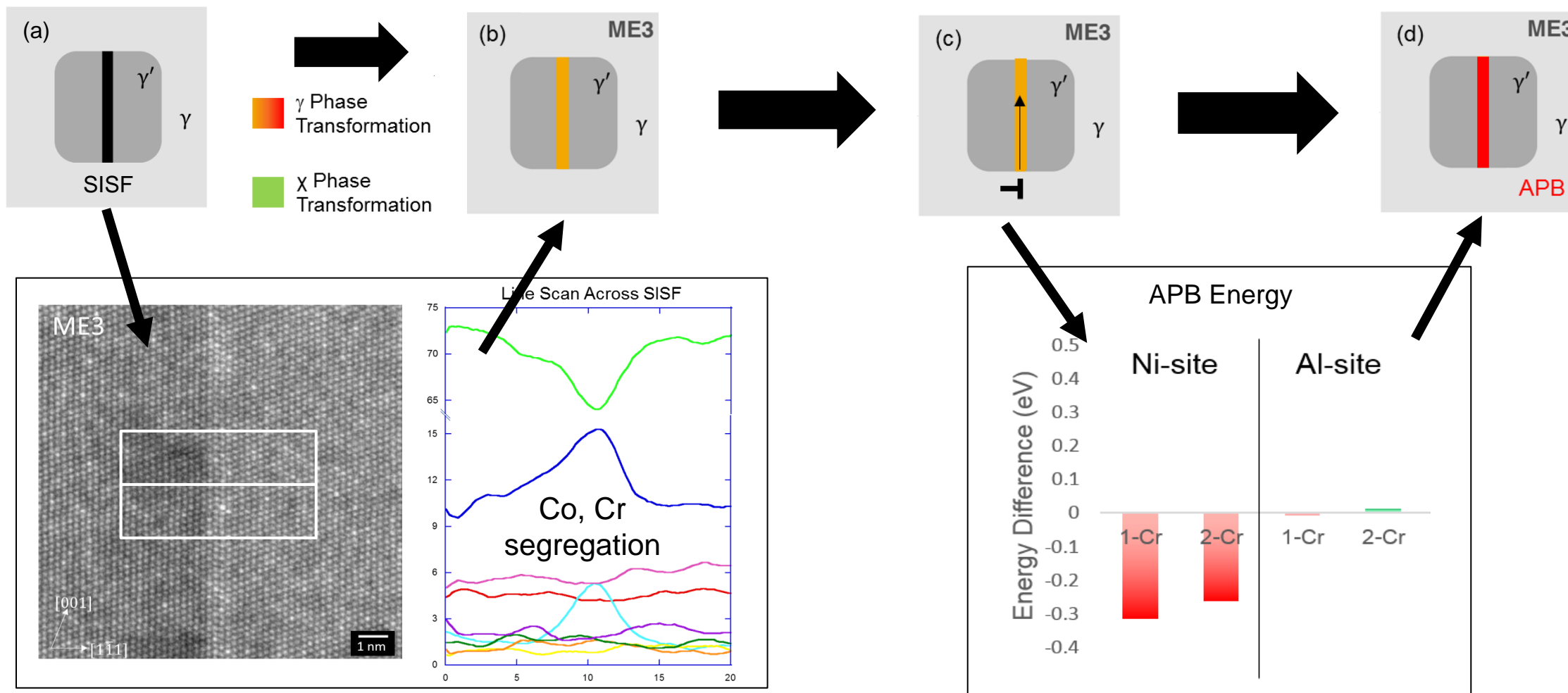
# Creep Performance of ME3 and LSHR



- Creep tests were performed at 760°C under a stress of 552MPa
- LSHR has consistently performed better in creep compared to ME3 in this temperature regime. Why?



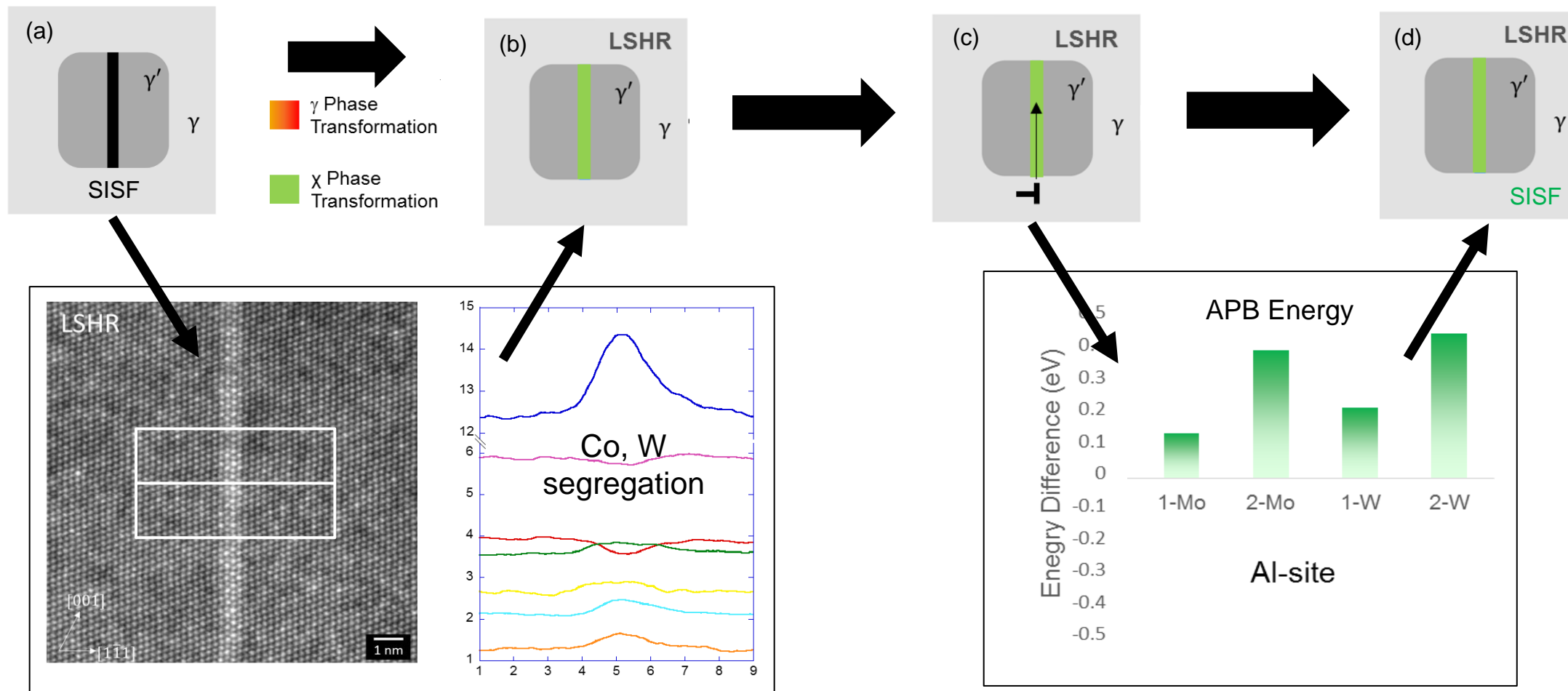
# Phase Transformation Softening – $\gamma$ Phase



$\gamma$  phase formation along SISF promotes stacking fault ribbon shear

SISF = Superlattice Intrinsic stacking Fault

# Phase Transformation Strengthening – $\chi$ Phase



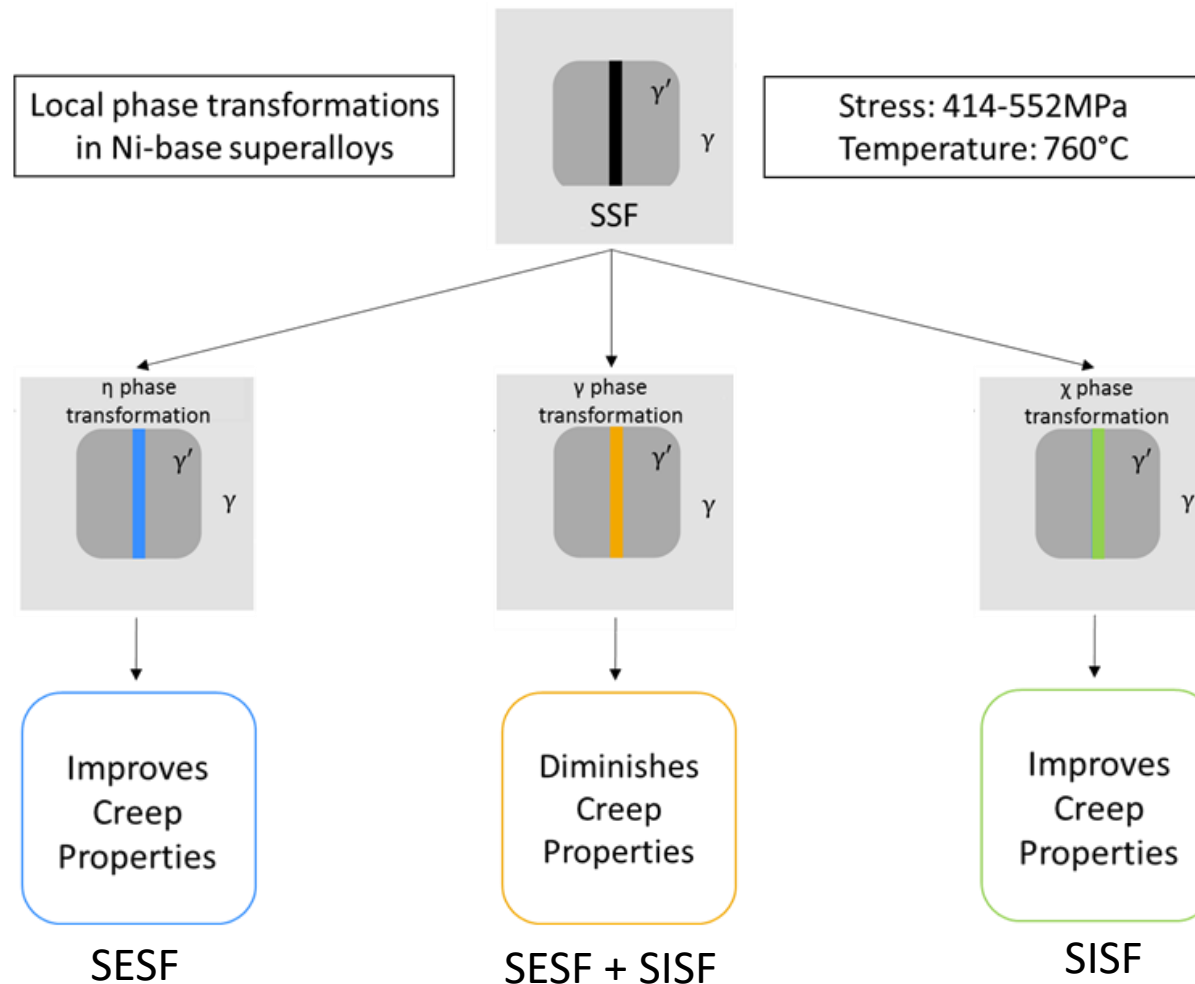
$\chi$  phase formation along SISF inhibits stacking fault ribbon shear

SISF = Superlattice Intrinsic stacking Fault





# Phase Transformation Strengthened Superalloys

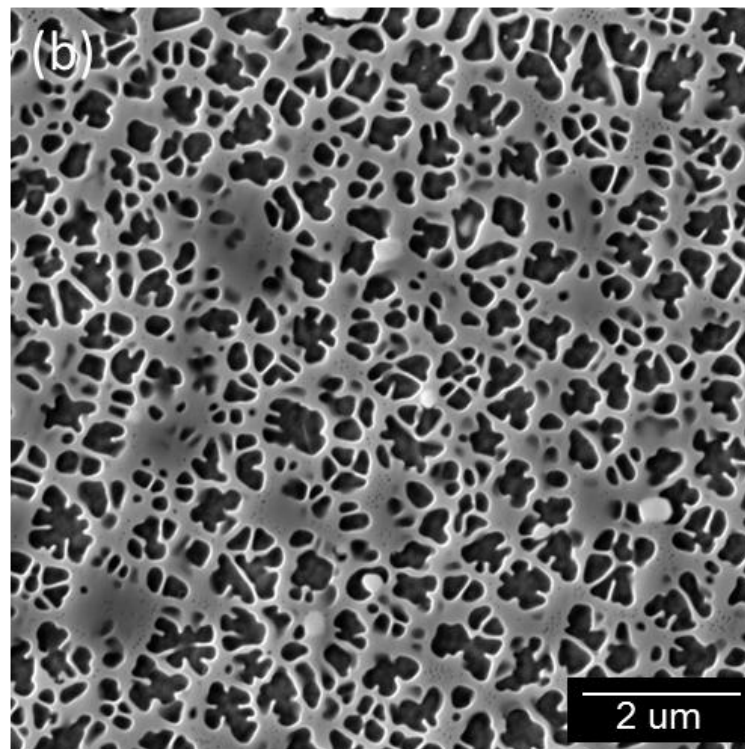
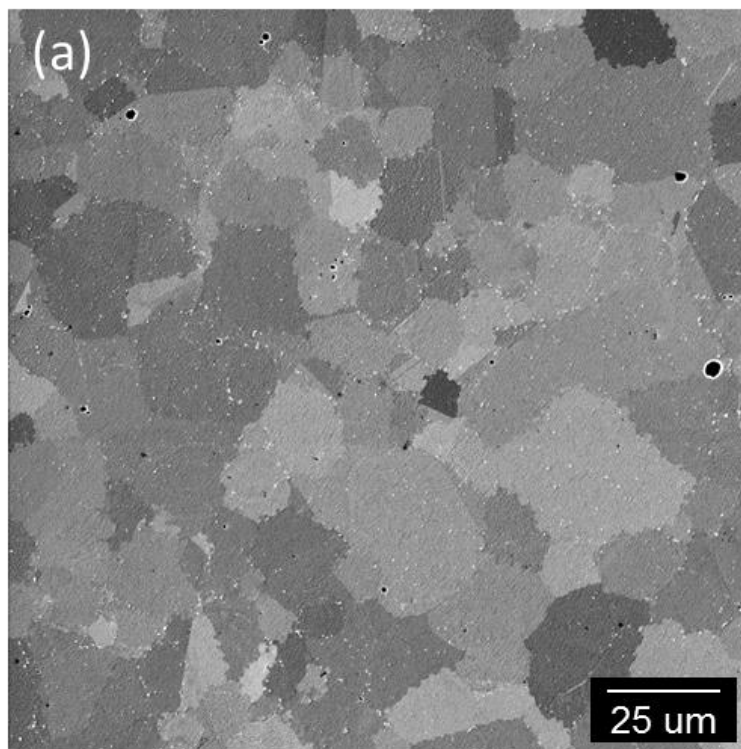


Can the  $\eta$  and  $\chi$  phase transformation strengthening mechanisms be combined into a single alloy without precipitating bulk topologically close packed (TCP) phases?



# Development of Transformation Strengthened NASA Alloys (TSNA)

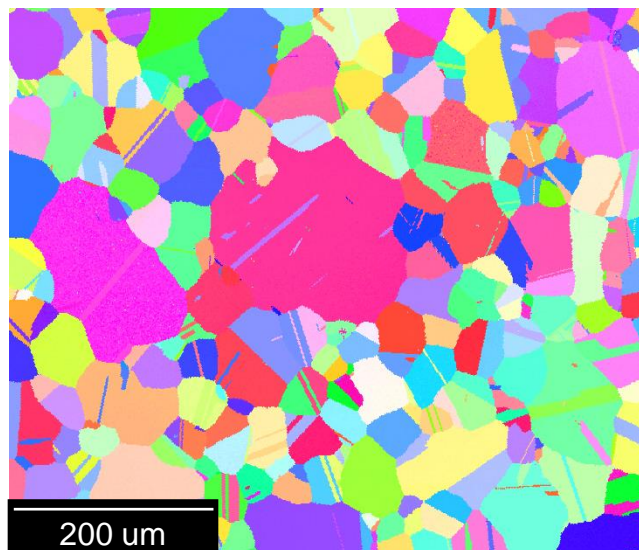
Alloy	Cr	Co	Al	Ti	Nb	Mo	Ta	W	Hf	B	C	Ni
LSHR	12.5	20.4	3.5	3.5	1.5	2.7	1.5	4.3	0	0.03	0.045	Bal
ME3	13	21	3.4	3.8	0.8	3.7	2.4	2.1	0	0.02	0.05	Bal
TSNA-1	10.9	19	2.9	3	1.4	2.6	5.0	4.5	0.37	0.025	0.05	Bal





# As-HIPed TSNA-1

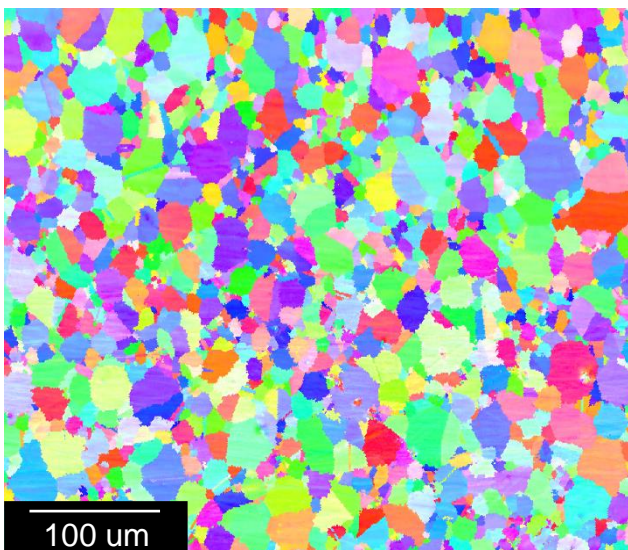
## LSHR - CG



200 um

Avg. Grain Size: 60um

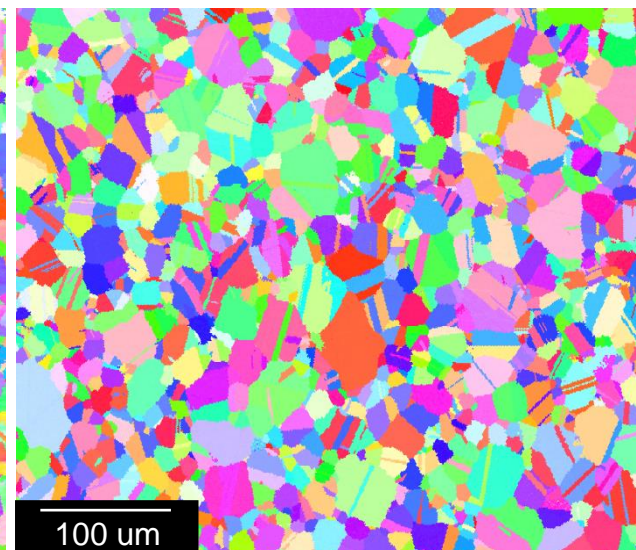
## TSNA-1 - FG



100 um

Avg. Grain Size: 19.2um

## LSHR - FG



100 um

Avg. Grain Size: 19.8um

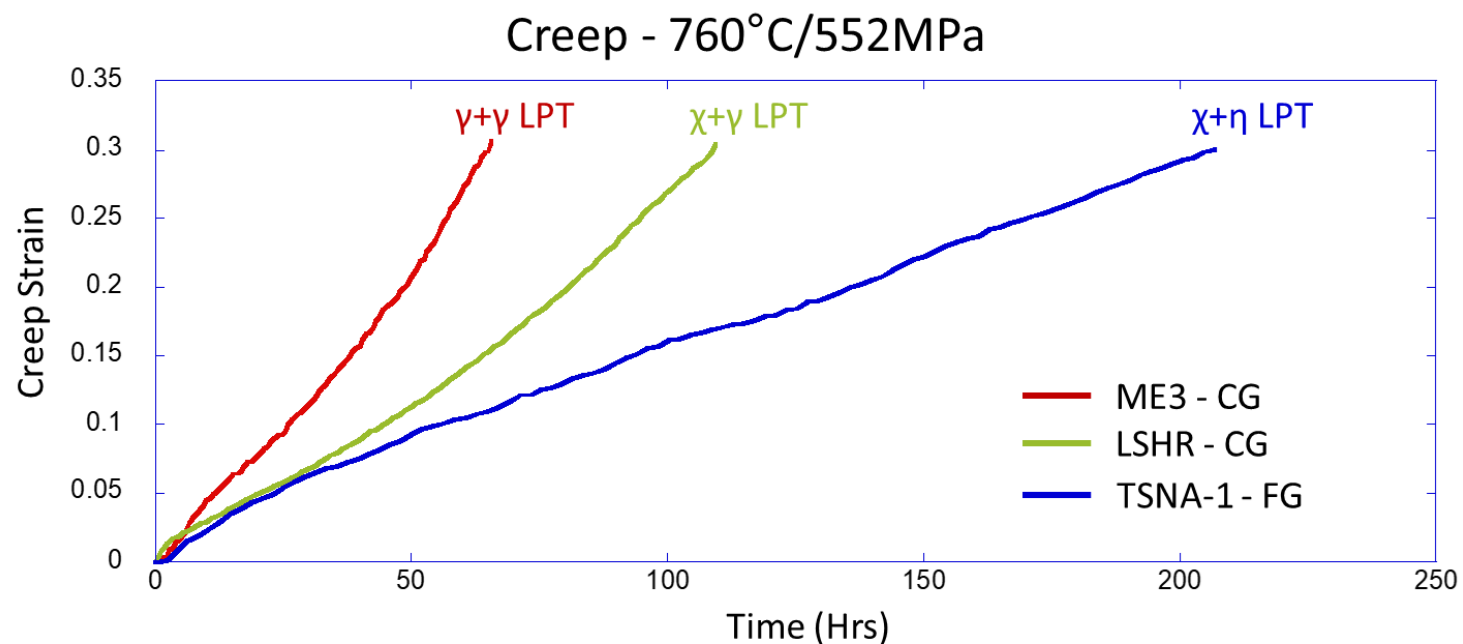
Alloy	Secondary $\gamma'$ VF	Tertiary $\gamma'$ VF	Total $\gamma'$ VF	Average Secondary $\gamma'$ Size	Average Tertiary $\gamma'$ Size
<b>ME3</b>	44.8 ± 0.5 %	2.6 ± 0.2 %	<b>47.4 ± 0.6 %</b>	234 nm	36.4 nm
<b>LSHR</b>	45.4 ± 1.8 %	3.4 ± 0.4 %	<b>48.7 ± 1.3 %</b>	243 nm	39.8 nm
<b>TSNA-1</b>	54.0 ± 0.2 %	0.6 ± 0.2 %	<b>54.5 ± 0.3 %</b>	311 nm	38.5 nm

By not forging the TSNA-1 alloy, grain sizes remained fine in comparison to LSHR. A fine grain LSHR was produced for a better comparison.





# Creep Properties

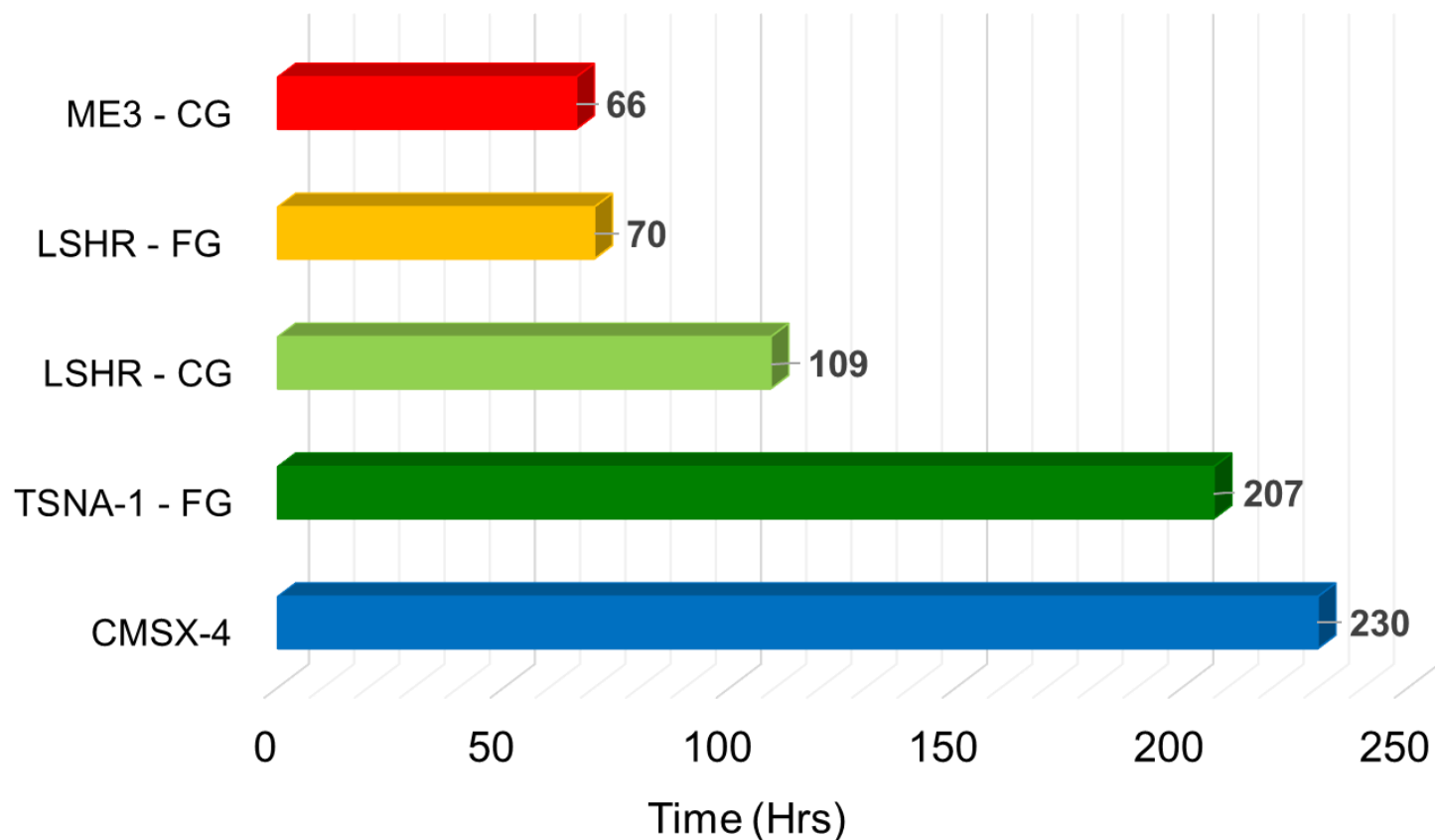


TSNA-1 presents significantly better creep properties over current state of the art alloys through possible phase transformation strengthening



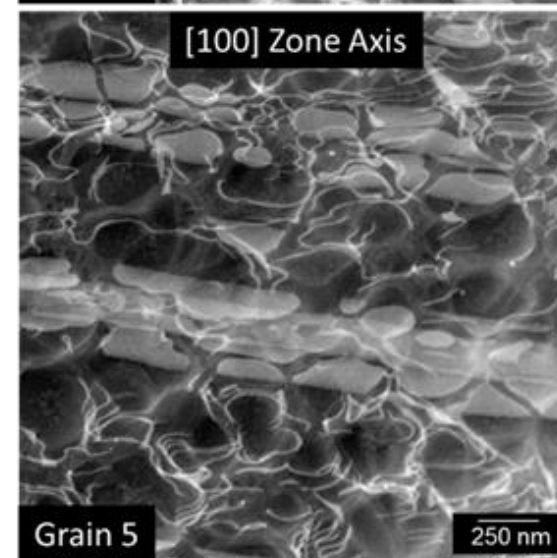
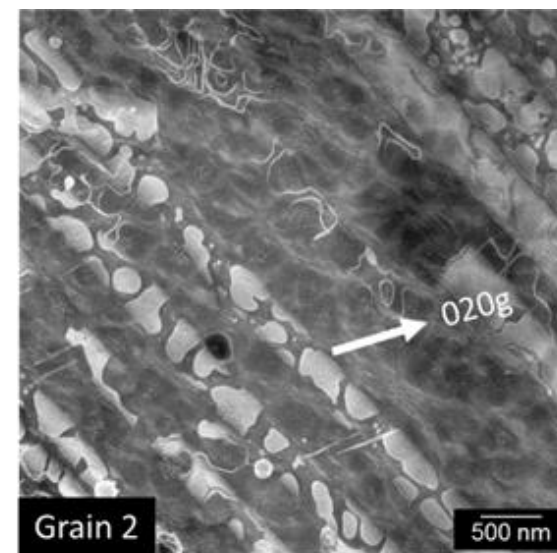
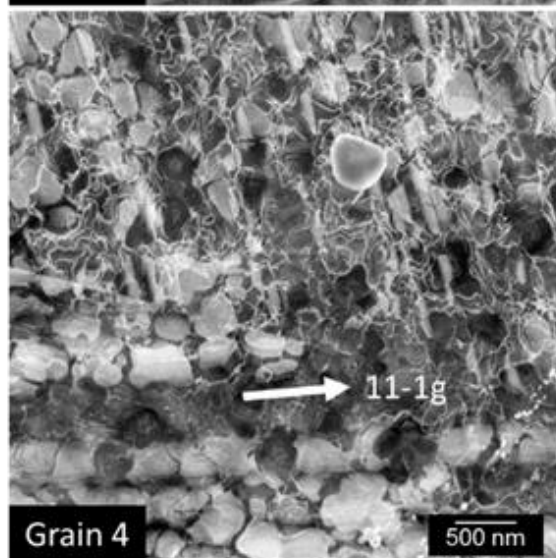
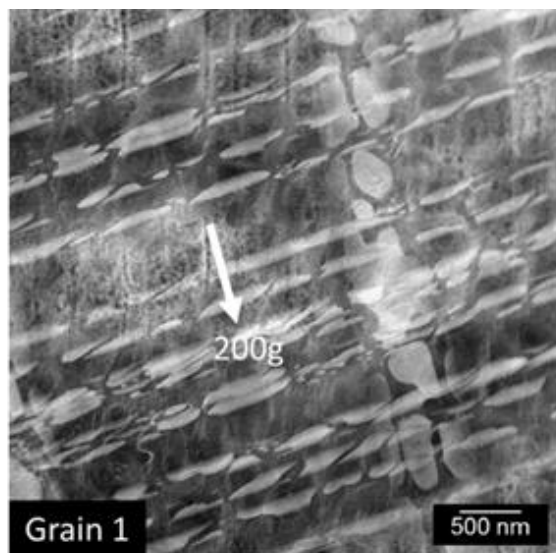
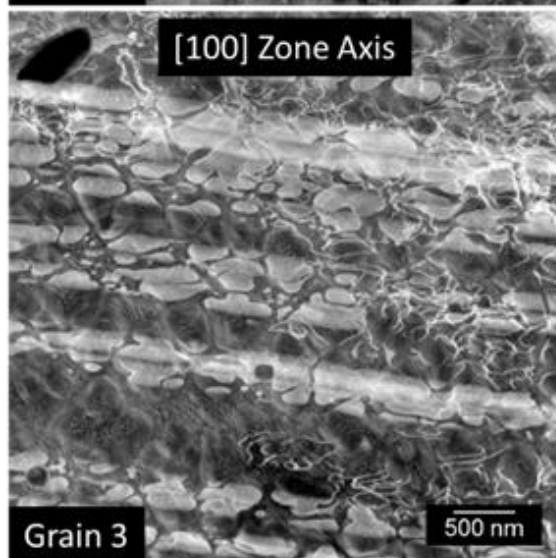
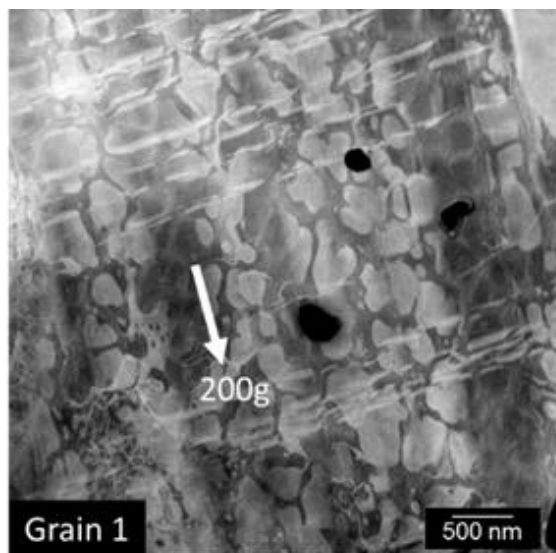
# Creep Properties

Time to 0.3% Creep Strain



TSNA-1 presents significantly better creep properties over current state of the art alloys through possible phase transformation strengthening

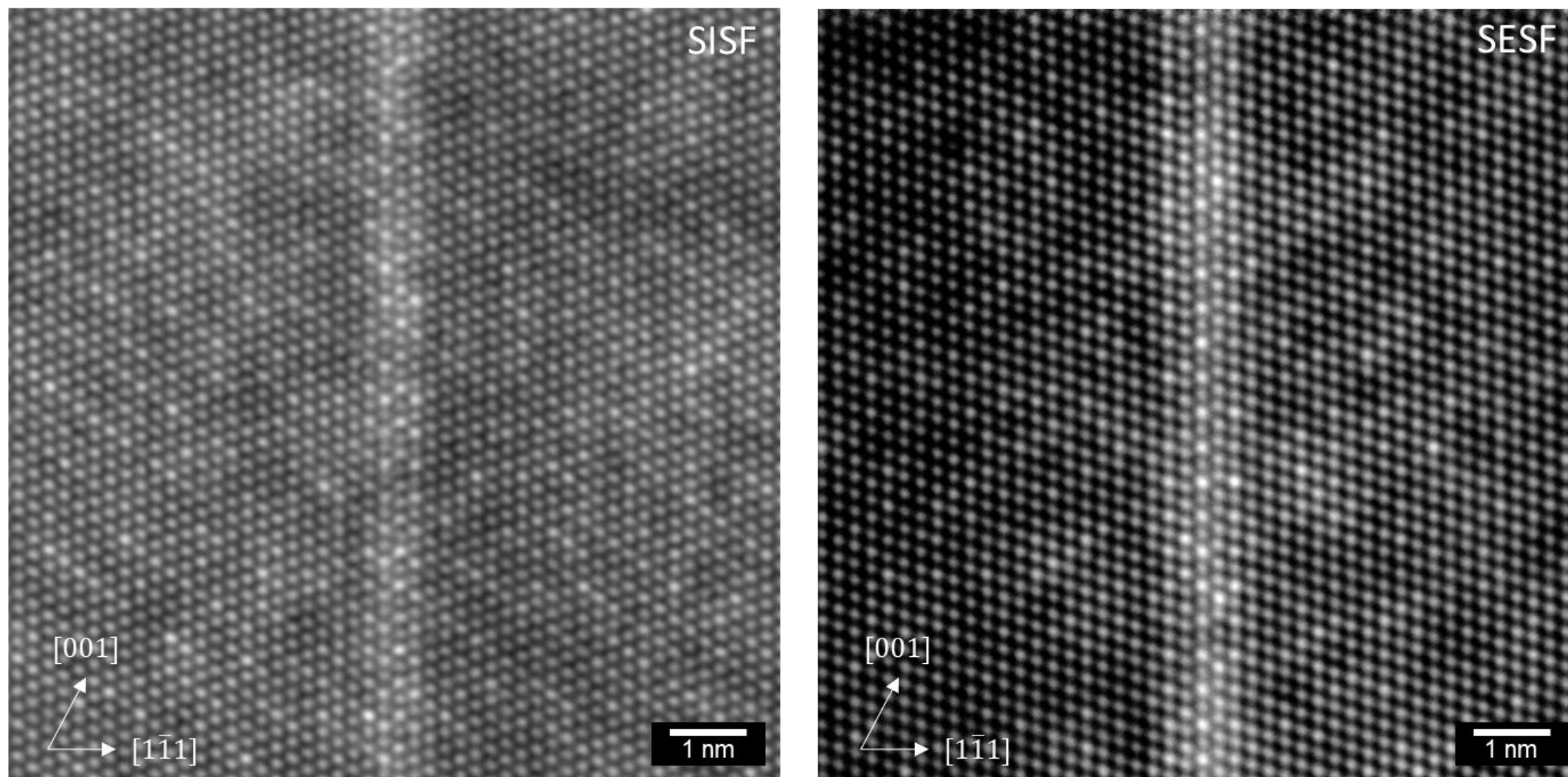
# Creep Deformation in TSNA-1



Deformation was dominated by isolated faulting in the  $\gamma'$  precipitates and dislocations gliding in the matrix.



# Stacking Fault Segregation in TSNA-1

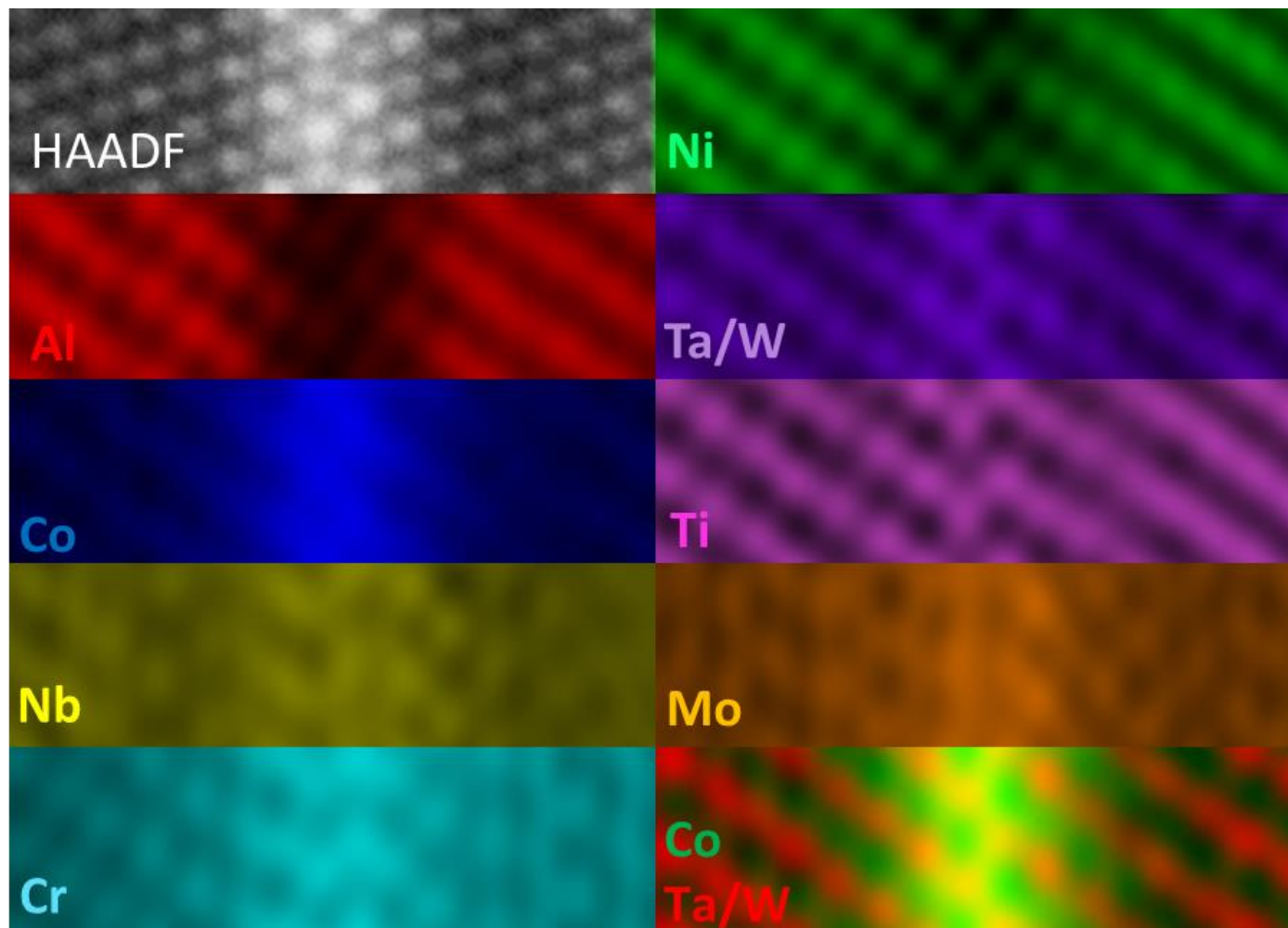
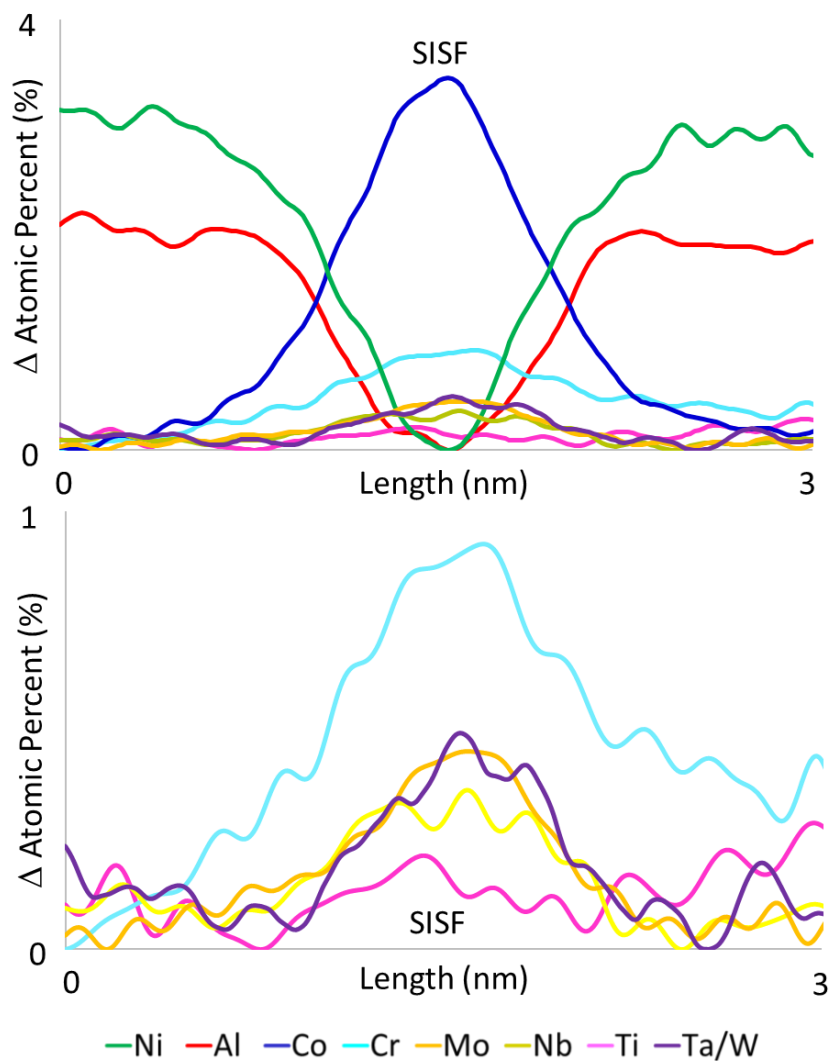


Ordered segregation observed in HAADF images of both fault types





# Phase Transformations in TSNA-1 - SISF



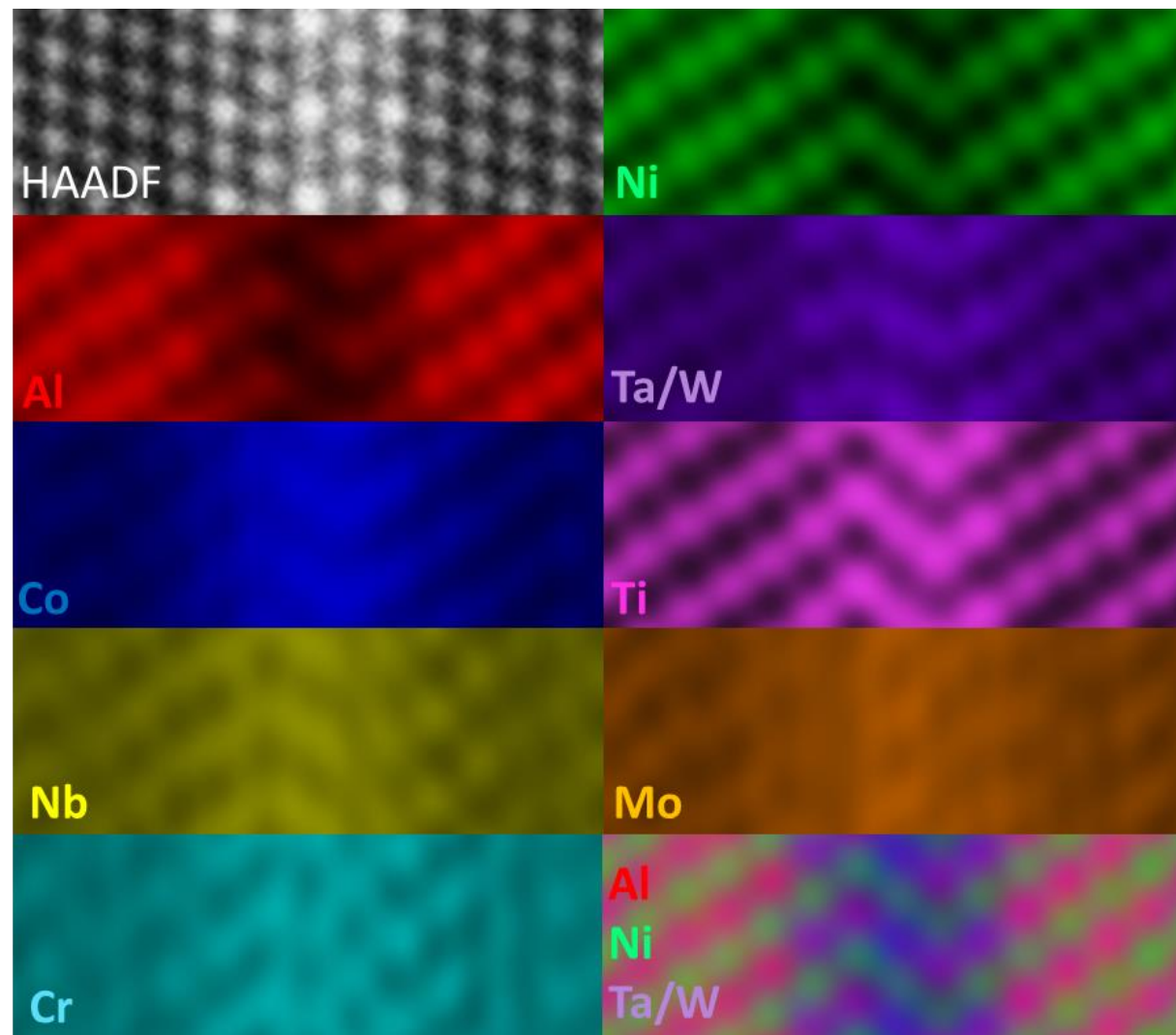
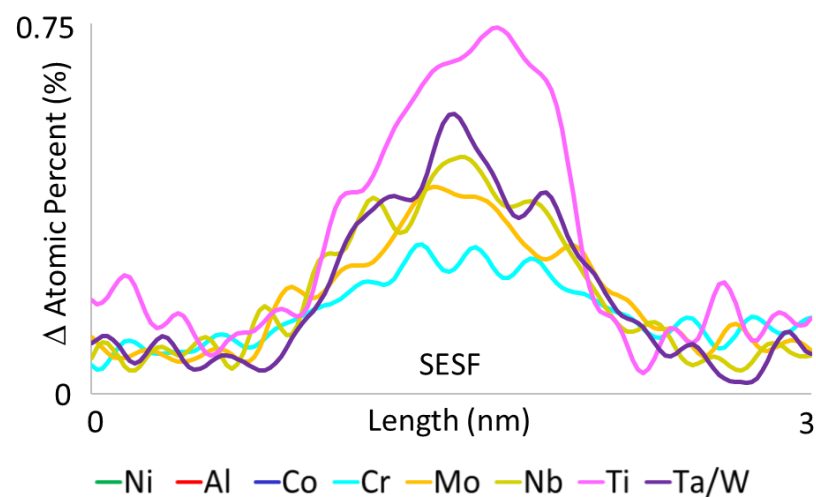
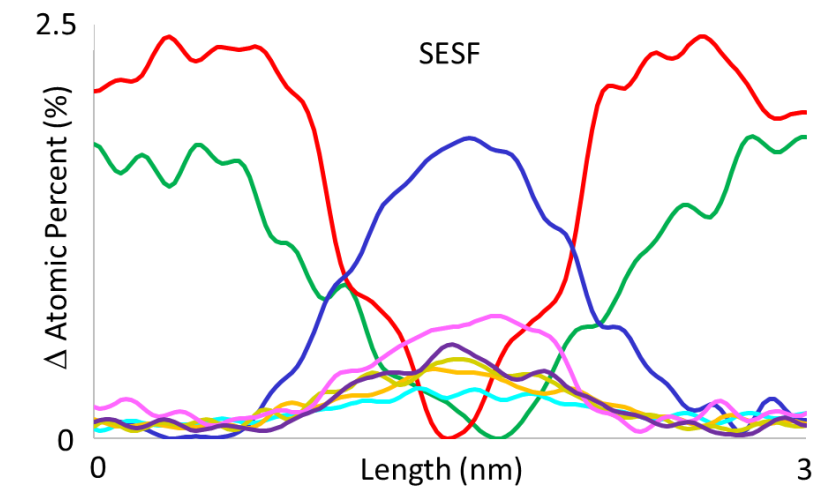
$\gamma'$  to  $\chi$  phase transformation confirmed along SISFs





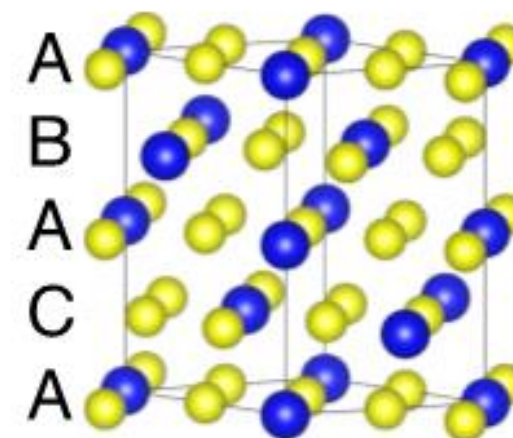
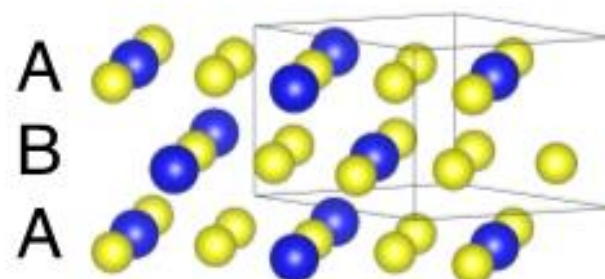
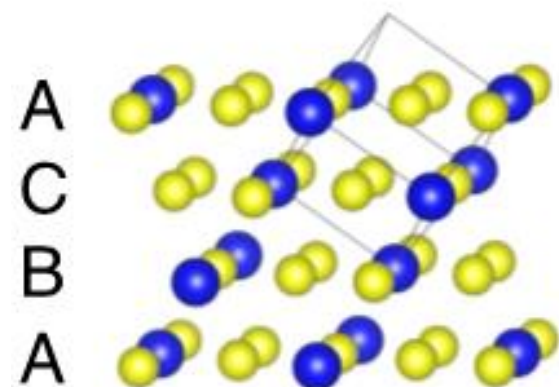
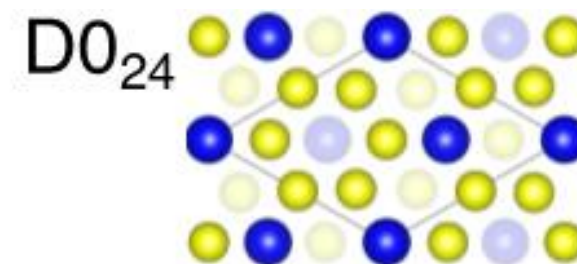
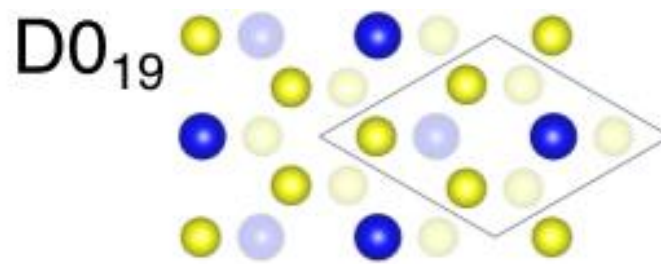
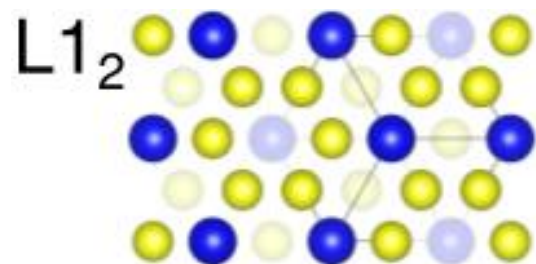


# Phase Transformations in TSNA-1 - SESF



$\gamma'$  to  $\eta$  phase transformation confirmed along SESFs

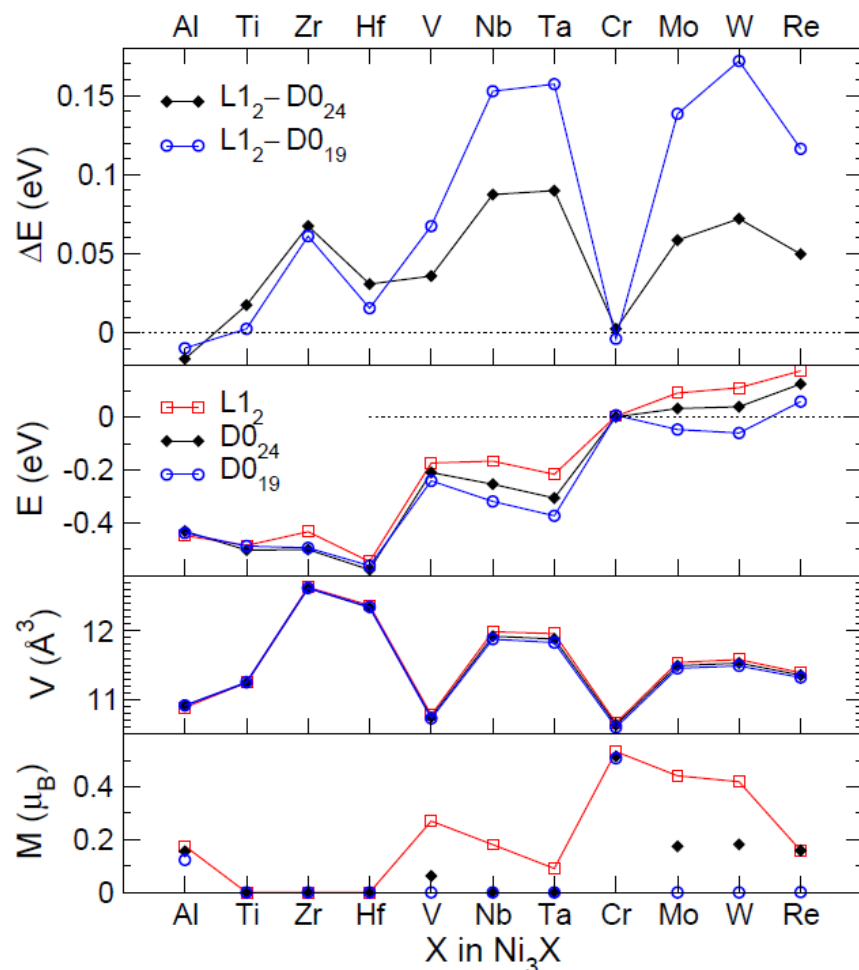
# Density Functional Theory Calculations



Binary  $L1_2$ ,  $D0_{19}$ , and  $D0_{24}$  cells were produced for energy calculations

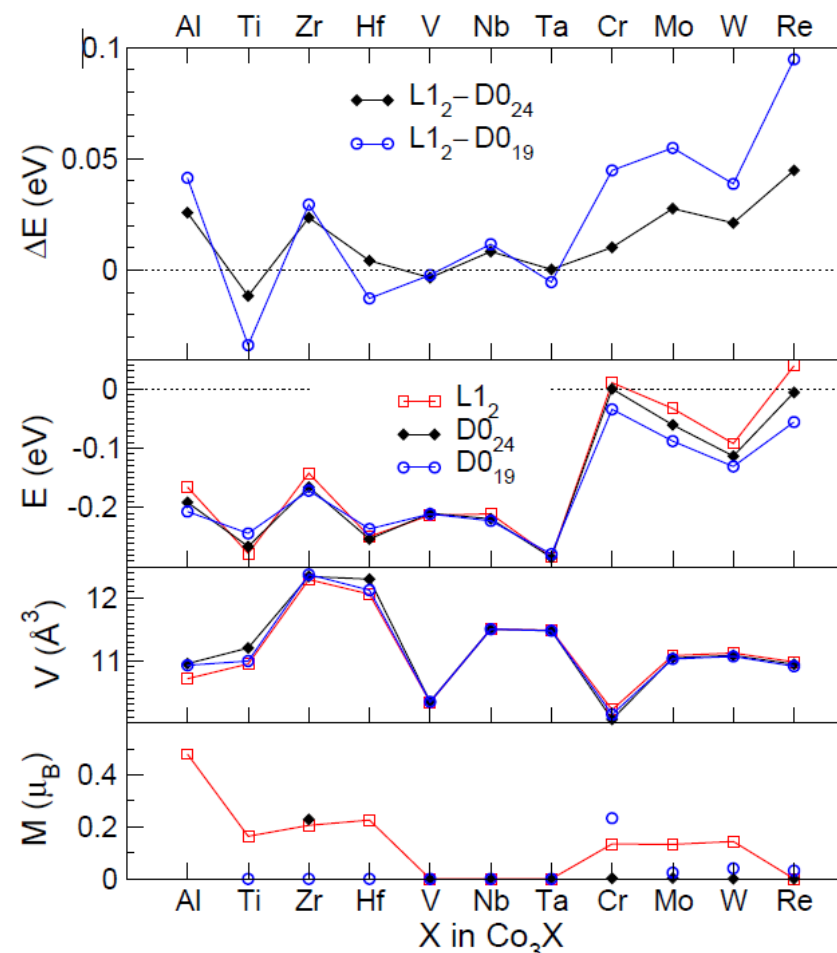
# Density Functional Theory Calculations

## Ni<sub>3</sub>X



$\eta$  formers: **Ta, Nb, W, Zr, V, Hf, Ti**  
 $\chi$  formers: **W, Ta, Nb, Mo, V, Zr**

## Co<sub>3</sub>X

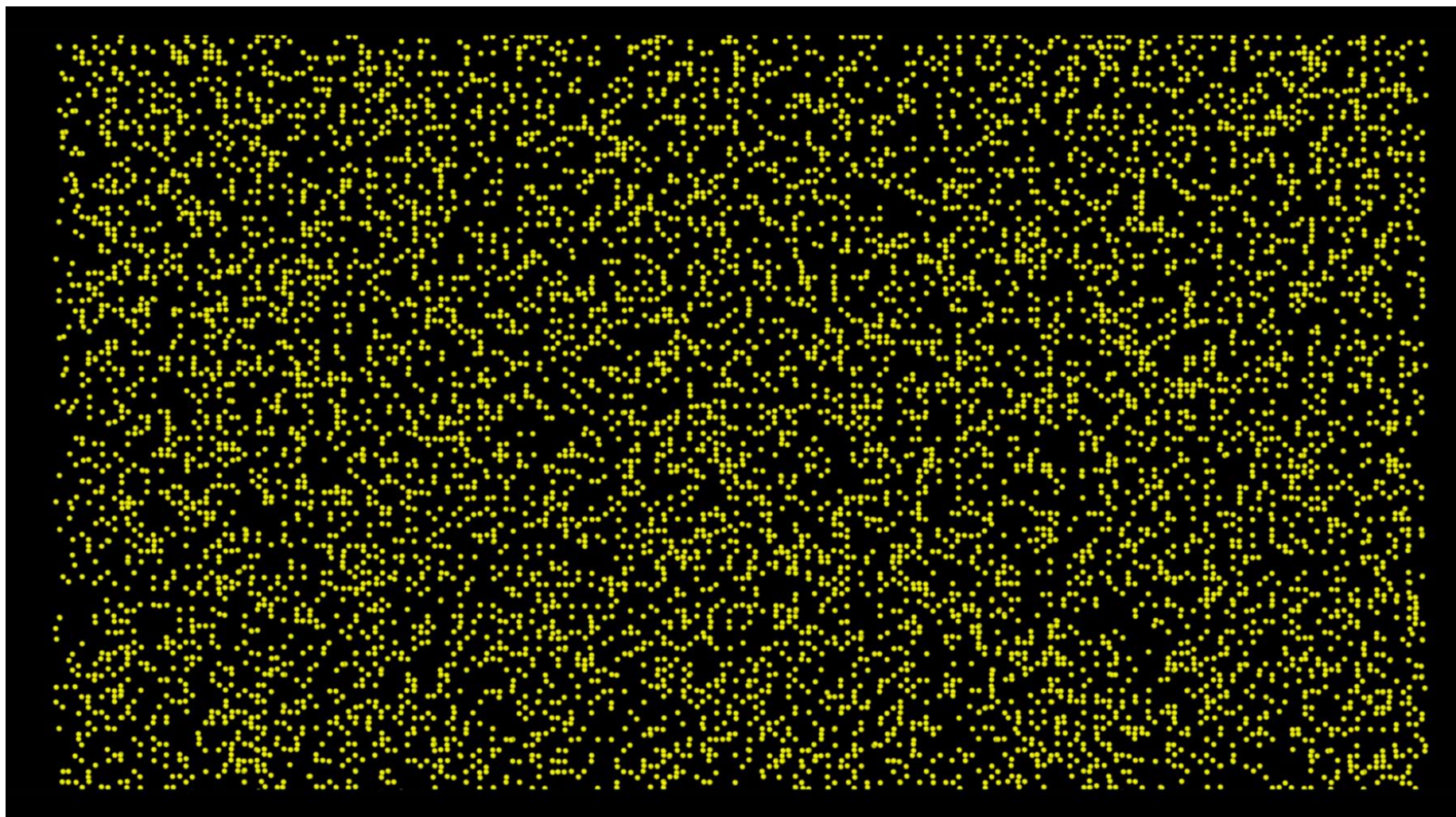


$\eta$  formers: **Re, Mo, W, Zr, Nb**  
 $\chi$  formers: **Re, Mo, Cr, W, Zr, Nb**

**Nb and Ta appear to be stronger promoters of  $\eta$  phase compared to Ti**



# Future Work: Molecular Dynamic Models – Nb Segregation along SISFs



$T = 1000 \text{ K}$

Ni and Al atoms not shown

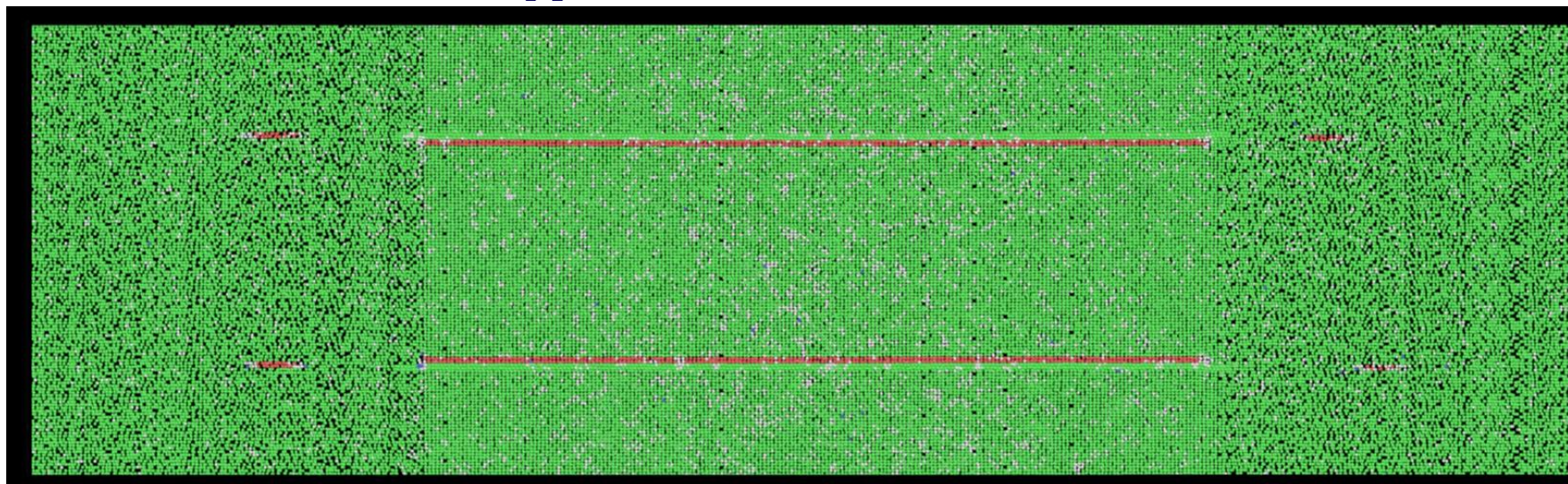
Substituted 10% Al in precipitate with Nb





# Future Work: Molecular Dynamic Models

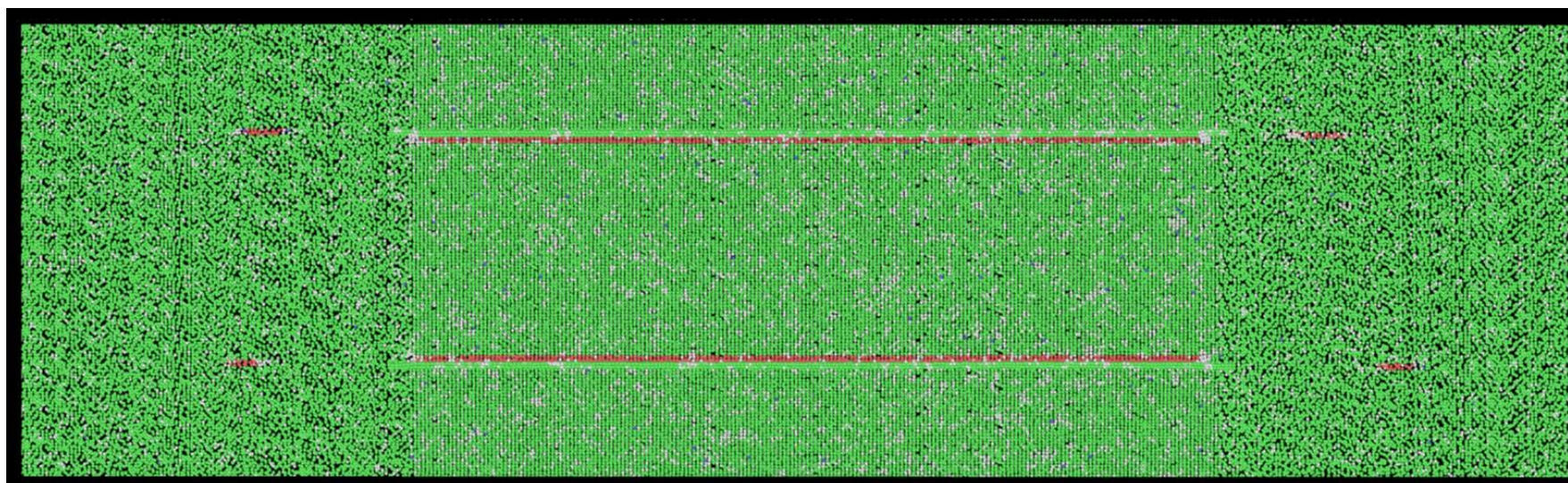
## - $\chi$ Phase Effect on Dislocation Motion



sxz = 400 MPa  
syz = 693 MPa  
stot = 800 MPa  
T = 1000 K

Ni atoms not shown

**0% Nb**



sxz = 400 MPa  
syz = 693 MPa  
stot = 800 MPa  
T = 1000 K

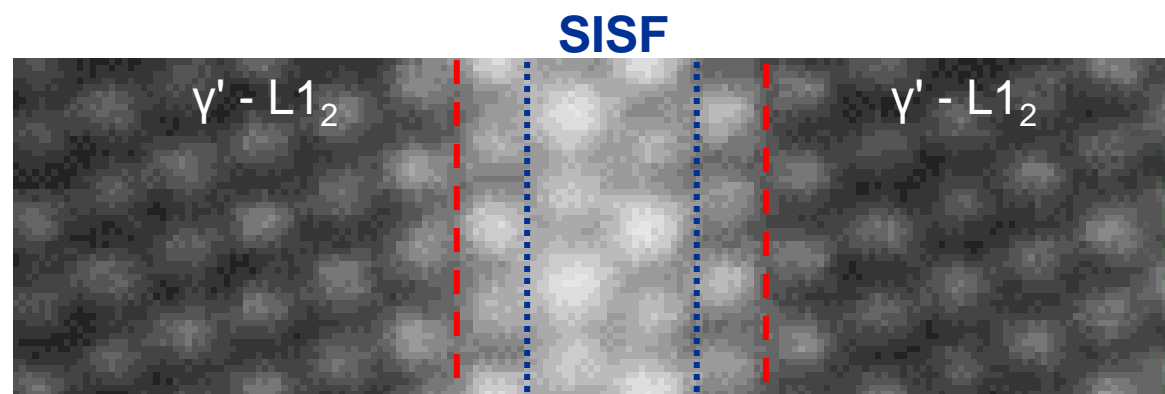
Ni atoms not shown

**7.5% Nb**

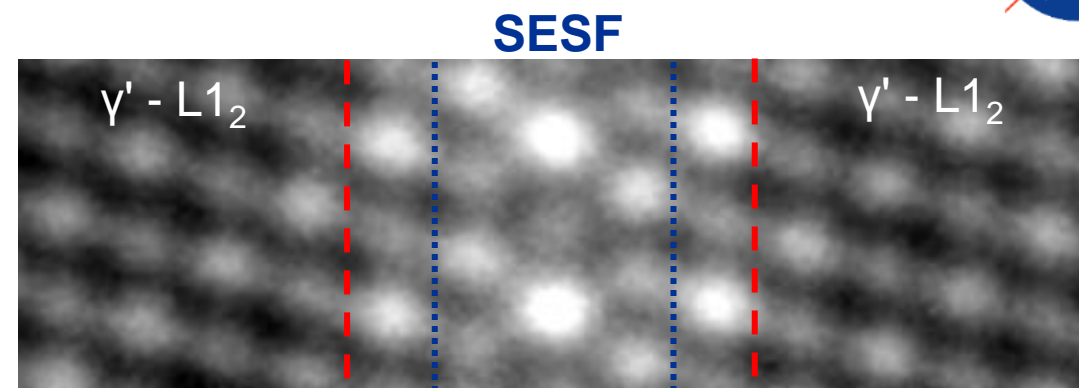
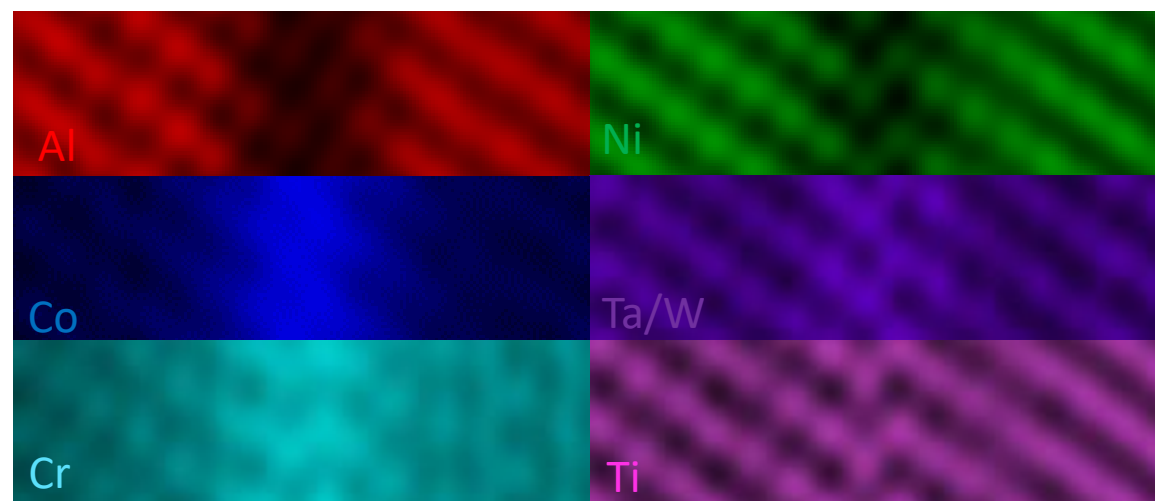




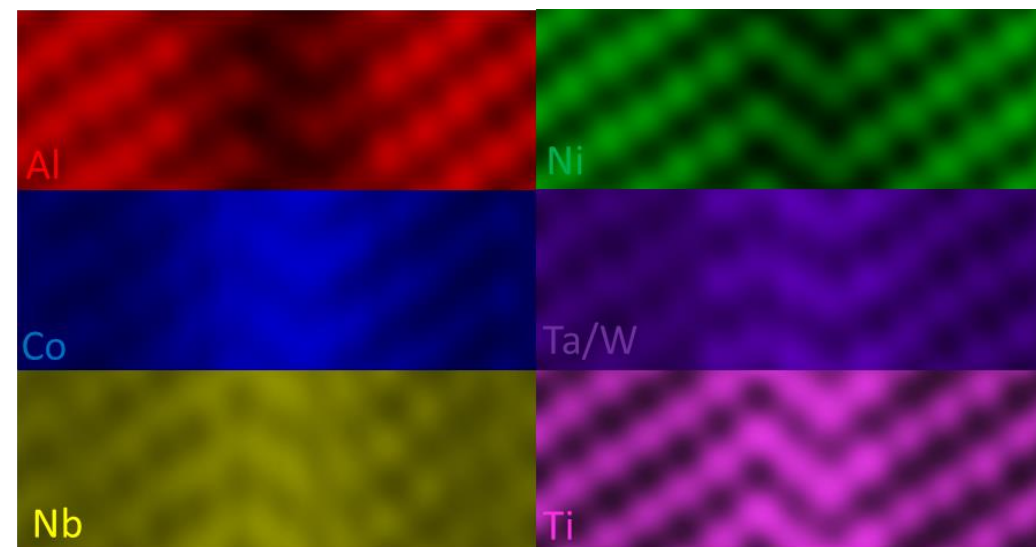
# Why Phase Transformations?



**D0<sub>19</sub>**



**D0<sub>24</sub>**



Local crystal structure + composition + observed Z contrast ordering = Local phase transformation along faults



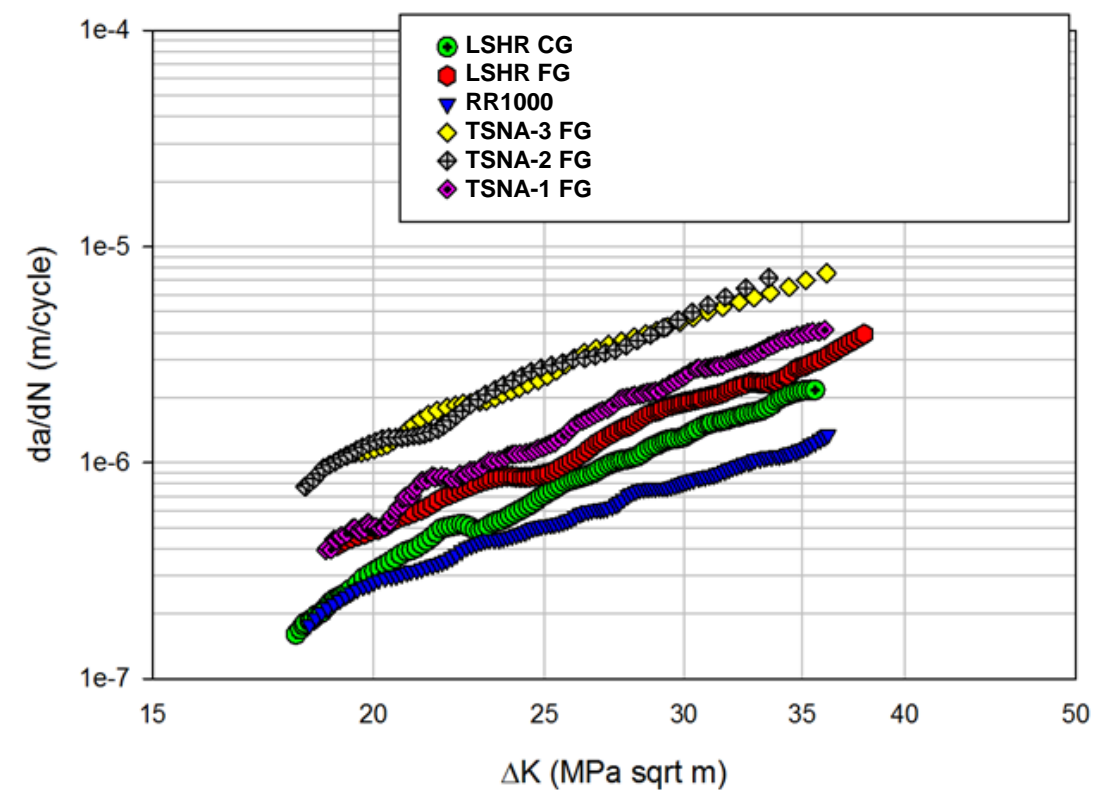


# Future Work

## Forged TSNA-1



## Crack Growth





# Conclusions

- The creep performance of TSNA-1 is significantly better compared to LSHR and ME3 despite testing conducted on an overall similar, though not yet optimized, microstructure.
- The creep deformation at 760°C/552MPa in TSNA-1 is dominated by dislocation glide in the  $\gamma$  channels and isolated faulting in the  $\gamma'$  precipitates.
- High resolution STEM analysis reveals the formation of  $\chi$  phase along SISFs and  $\eta$  phase along SESFs for TSNA-1.
- The formation of these phases along the faults may explain the superior creep properties exhibited by TSNA-1, as the grain and  $\gamma'$  microstructure fail to do so.
- **The strengthening  $\eta$  and  $\chi$  phase transformations can be combined in future Ni-base disk alloy compositions for improved creep properties**





# Acknowledgements

## Questions?



- CEMAS (OSU)
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- Bob Carter
- Cheryl Bowman
- Rick Rogers
- Steve Arnold
- Jack Telesman
- Pete Bonacuse
- Joy Buehler
- Alex Leary

